

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

REGION 5 CHICAGO REGIONAL LABORATORY

536 SOUTH CLARK STREET





Date: 1/13/2015

Subject: Review of Region 5 Data for BP Whiting Refinery

To: Air Division, US EPA Region 5

77 West Jackson Boulevard

Chicago, IL 60605

From: Wayne Whipple, Analyst

US EPA Region 5 Chicago Regional Laboratory

The data transmitted under this cover memo successfully passed CRL's data review procedures as documented in the current Quality Management Plan and applicable Standard Operating Procedures. In accordance with EPA's *Guidance on Environmental Data Verification and Data Validation* (Document EPA QA/G-8), CRL verified and validated the data but does not perform data quality assessment based on project plans.

This report was reviewed and the information provided herein accurately represents the analysis performed.

Jan 13 2015

Please contact the analyst with any technical report issues, Amanda Wroble at (312)-353-0375 for sample project concerns, and Sylvia Griffin at (312)-353-9073 with data transmittal questions. Thank you.

Attached are Results for: BP Whiting Refinery

Data Management Coordinator and DateTransmitted

Analyses included in this report:

Air Toxics Reimer 5

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Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605 Phone:(312)353-8370 Fax:(312)886-2591

Air Division, US EPA Region 5 Project: BP Whiting Refinery
77 West Jackson Boulevard Project Number: [none]

Chicago IL, 60605 Project Manager: Motria Caudill

Reported: Jan-13-15 14:58

Analysis Case Narrative for Volatile Organic Compounds (VOC) Air Toxics

Wayne J. Whipple, Ph.D. phone (312) 353-9063 email whipple.wayne@epa.gov

General Information

Three 15 L canisters and two six liter canisters were received in good condition from the Air and Radiation Division monitoring group on September 18, 2014. The samples met hold time.

The samples were analyzed for VOC Air Toxics using SOP MS-005 Revision 6 with cold trap dehydration preconcentrator on Amelia. (Reference Method US EPA TO-15). Three Pen and Ink Changes are also included, MS005 r5.0-P01 and P03. The pen and ink changes use the manufacturers recommended mass spectrometer tune check requirements instead of BFB. The halocarbon sample monitoring compound concentrations and acceptance criteria is updated in -03.

Data are reported to a reporting limit of 25 pptv and at least 35 pptv was requested in the Quality Assurance Project Plan, QAPP and meets the QAPP requirements. (EPA Region 5 VOC Sampling Comparison QAPP v1 1)

Sample Analysis and Results

The criteria used for determining if system contamination influences the sample results is for a sample result to be over three to five times the contamination in the blank checks.

Compounds reported have been updated since MS 005 revision 6. The standards that are available do not contain many analytes that are listed in the SOP. There are a few compounds in the standard that have not been fully documented for CRL quality assurance. Since they have not undergone full quality assurance acceptance these compounds, n-heptane and naphthalene, 4-ethyl toluene, *trans* 1,2-dichloroethene, and dibromochloromethane will be flagged as estimated.

Quality Control

Calibration modifications:

The sample results for acetone in all the samples were higher than the highest calibration curve for sample 1409007-02 which was not run for dilution so the data was flagged estimated in that sample. Other sample results



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for acetone and ethanol were reported from diluted analysis. The QC results for those samples are only presented for the reported analytes and selected sample monitoring compounds.

Data from 1409007-04 and -05 have been reanalyzed multiple times at various dilutions. Acetone, Methylene Chloride and isopropyl alcohol are rejected from sample 1409007-04 and is reported from the reanalysis 1409007-04RE2. The QC from that batch is only reported for those two compounds.

Acrolein and 1,3-butadiene and propene are biased high because of coeluting hydrocarbons that cannot be resolved or deconvoluted and therefore the data is flagged biased high, K..

A number of compounds had their reporting limits raised to an acceptable region of the calibration curve that passes the SOP criteria. Ethanol, 1,2-dichlorobenzene, o-xylene, and isopropyl alcohol are raised to 50 pptv. Methylene chloride, styrene and bromoform, m+p-xylene, 1,3- and 1,4-dichlorobenzene and hexachlorobutadiene are raised to 100 pptv and 4-methyl-2-pentanone rejected because of vary poor calibration performance.

There are two calibration file date and time stamps because of the software updates the date and time stamp for an unknown reason but the equations have not changed. The two reports are provided to show identical equations and calculations.

F-113 and 1,1-dichloroethene had low recoveries in the Initial Calibration Verification (ICV) standard and are therefore flagged as estimated.

1,4-dioxane's and 2-hexanone's calibration had an excessive percent error that does not meet the SOP criteria throughout most of the calibration curve and therefore the results are rejected.

Relative percent difference, RPD results are followed from the SOP and not the final report.

Ethanol is qualified as estimated "J" flag because of poor calibration performance. The calibration is not ideal and reduces the confidence in the results for this compound.

Naphthalene is flagged as estimated because the performance of the analytical system is not as reliable having an acceptable calibration response above 250 pptv for that late eluting compound and it fails the second source ICV standard.

1,1,1-Trichloroethane is reported below the detection limit for tracking of the sample monitoring compound. The result is within the expected global results, although is still flagged as estimated because it is below the detection limit. The result is also within 5 times the blank result but it is already flagged as estimated so no further action is required.

Carbon disulfide's analytical performance is extremely poor so the compound is rejected from this report.

Method Blank and Canister Blanks are qualified as high bias or have their reporting limits raised if the results



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from the sample are within 5x the results from the blank. Acetone has been flagged K as high bias for the diluted samples although the result is very close to the undiluted sample suggesting the bias may be negligible. Methylene chloride is qualified high bias in samples -04 and -05.

System monitoring compounds are guidelines to monitor the success and quality of analytical integrity. They are guidelines with limits set around very consistent global background concentrations. The limits are suggested and open to interpretation of the analyst. Many of the compounds are extremely close to or below the detection limit of the instrumentation. There is no reason to flag any of the samples in this data set for exceedences of the limits applied to these compounds.

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ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received	
4	1409007-01	Air	Sep-18-14 09:39	Sep-18-14 14:00	
1	1409007-02	Air	Sep-18-14 10:25	Sep-18-14 14:00	
3	1409007-03	Air	Sep-18-14 11:20	Sep-18-14 14:00	
2	1409007-04	Air	Sep-18-14 12:10	Sep-18-14 14:00	
4B	1409007-05	Air	Sep-18-14 00:00	Sep-18-14 14:00	

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Project: BP Whiting Refinery

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Air Toxics by GC/MS, EPA TO-15 (modified) US EPA Region 5 Chicago Regional Laboratory

4 (1409007-01) Air Sampled: Sep-18-14 09:39 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Propene	1.77	K		0.0250	ppbv	1	B409106	Sep-19-14	Sep-19-14
Chloromethane	0.496			0.0250	"	"	"	"	"
Vinyl chloride	U			0.0250	"	"	"	"	"
1,3-butadiene	0.0440	K		0.0250	"	"	"	"	"
Bromomethane	U			0.0250	"	"	"	"	"
Chloroethane	U			0.0250	"	"	"	"	"
Acrolein	0.891	K		0.0250	"	"	"	"	"
Isopropyl alcohol	1.52	J		0.0500	"	"	"	"	"
1,1-Dichloroethene	U	J		0.0250	"	"	"	"	"
Methylene chloride	0.259			0.100	"	"	"	"	"
Carbon disulfide	Rejected			0.0250	"	"	"	"	"
Methyl tert-butyl ether	U			0.0250	"	"	"	"	"
1,1-Dichloroethane	U			0.0250	"	"	"	"	"
Vinyl acetate	U			0.0250	"	"	"	"	"
Hexane	0.613			0.0250	"	"	"	"	"
2-Butanone	0.593			0.0250	"	"	"	"	"
cis-1,2-Dichloroethene	U			0.0250	"	"	"	"	"
trans-1,2-Dichloroethene	U	J		0.0250	"	"	"	"	"
Chloroform	0.0311			0.0250	"	"	"	"	"
1,2-Dichloroethane	U			0.0250	"	"	"	"	"
Cyclohexane	0.184			0.0250	"	"	"	"	"
Benzene	0.318			0.0250	"	"	"	"	"
n-Heptane	0.177	J		0.0250	"	"	"	"	"
1,2-Dichloropropane	U			0.0250	"	"	"	"	"
Trichloroethene	U			0.0250	"	"	"	"	"
Bromodichloromethane	U			0.0250	"	"	"	"	"
1,4-Dioxane	Rejected			0.0250	"	"	"	"	"
cis-1,3-Dichloropropene	U			0.0250	"	"	"	"	"
4-Methyl-2-pentanone	Rejected			0.0250	"	"	"	"	"
trans-1,3-Dichloropropene	U			0.0250	"	"	"	"	"
1,1,2-Trichloroethane	\mathbf{U}			0.0250	"	"	"	"	"
Dibromochloromethane	\mathbf{U}	J		0.0250	"	"	"	"	"
Toluene	0.639			0.0250	"	"	"	"	"
2-Hexanone	Rejected			0.0250	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			0.0250	"	"	"	"	"
Tetrachloroethene	U			0.0250	"	"	"	"	"



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4 (1409007-01) Air Sampled: Sep-18-14 09:39 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Chlorobenzene	U			0.0250	ppbv	1	B409106	Sep-19-14	Sep-19-14
Ethylbenzene	0.0810			0.0250	"	"	"	"	"
m+p-Xylene	0.245			0.100	"	"	"	"	"
Bromoform	U			0.100	"	"	"	"	"
Styrene	U			0.100	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			0.0250	"	"	"	"	"
o-Xylene	0.0822			0.0500	"	"	"	"	"
4-ethyltoluene	0.108	J		0.0250	"	"	"	"	"
1,3,5-Trimethylbenzene	U			0.0500	"	"	"	"	"
1,2,4-Trimethylbenzene	0.137			0.0250	"	"	"	"	"
1,3-Dichlorobenzene	U			0.100	"	"	"	"	"
Benzyl chloride	U			0.0250	"	"	"	"	"
1,4-Dichlorobenzene	U			0.100	"	"	"	"	"
1,2-Dichlorobenzene	U			0.0500	"	"	"	"	"
1,2,4-Trichlorobenzene	0.200			0.100	"	"	"	"	"
Naphthalene	1.22	J, K		0.100	"	"	"	"	"

Surogate	Result		%REC	%REC Limits	Batch	Prepared	Analyzed
Surrogate: Dichlorodifluoromethane	0.353		67%	90-120	"	"	"
Surrogate: Dichlorotetrafluoroethane	0.0189		95%	80-120	"	"	"
Surrogate: Trichlorofluoromethane	0.200		84%	90-120	"	"	"
Surrogate: 1,1,2-trichloro-1,2,2-trifluoroethane (Fi	0.0357	J	49%	90-120	"	"	"
Surrogate: 1,1,1-Trichloroethane	4.36E-3	J	87%	80-120	"	"	"
Surrogate: Carbon tetrachloride	0.0663		77%	80-120	"	"	"

4 (1409007-01RE1) Air Sampled: Sep-18-14 09:39 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Ethanol	2.40	J		0.500	ppbv	10	B411044	Sep-22-14	Sep-22-14
Acetone	3.55	J, K		0.250	"	"	"	"	"
Surogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Surrogata: Dichlorodifluoromathana	0.490			03%		00.120	"	"	"

Surrogate: Dichlorodifluoromethane 0.490 93% 90-120 " " "
Surrogate: Trichlorofluoromethane U 83% 90-120 " " "
Surrogate: Carbon tetrachloride U 67% 80-120 " " "



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Air Toxics by GC/MS, EPA TO-15 (modified) US EPA Region 5 Chicago Regional Laboratory

1 (1409007-02) Air Sampled: Sep-18-14 10:25 Received: Sep-18-14 14:00

	ъ .	Flags /	1001	Reporting	TT 1	D3:	D. I	ъ .	
Analyte	Result	Qualifiers K	MDL	Limit	Units	Dilution	Batch	Prepared	Analyzed
Propene		K		0.0250	ppbv "	1 "	B409106	Sep-19-14	Sep-19-14
Chloromethane	0.495			0.0250	"	"	"	"	"
Vinyl chloride	U			0.0250					
1,3-butadiene	0.0356	K		0.0250	"	"	"	"	"
Bromomethane	U			0.0250	"	"	"	"	"
Chloroethane	U			0.0250	"	"	"	"	"
Ethanol	1.94	J		0.0500	"	"	"	"	"
Acrolein	0.760	K		0.0250	"	"	"	"	"
Isopropyl alcohol	1.11	J		0.0500	"	"	"	"	"
Acetone	2.78	J		0.0250	"	"	"	"	"
1,1-Dichloroethene	U	J		0.0250	"	"	"	"	"
Methylene chloride	0.182			0.100	"	"	"	"	"
Carbon disulfide	Rejected			0.0250	"	"	"	"	"
Methyl tert-butyl ether	U			0.0250	"	"	"	"	"
1,1-Dichloroethane	U			0.0250	"	"	"	"	"
Vinyl acetate	U			0.0250	"	"	"	"	"
Hexane	0.543			0.0250	"	"	"	"	"
2-Butanone	0.353			0.0250	"	"	"	"	"
cis-1,2-Dichloroethene	U			0.0250	"	"	"	"	"
trans-1,2-Dichloroethene	U	J		0.0250	"	"	"	"	"
Chloroform	0.0316			0.0250	"	"	"	"	"
1,2-Dichloroethane	U			0.0250	"	"	"	"	"
Cyclohexane	0.237			0.0250	"	"	"	"	"
Benzene	0.283			0.0250	"	"	"	"	"
n-Heptane	0.227	J		0.0250	"	"	"	"	"
1,2-Dichloropropane	U			0.0250	"	"	"	"	"
Trichloroethene	U			0.0250	"	"	"	"	"
Bromodichloromethane	U			0.0250	"	"	"	"	"
1,4-Dioxane	Rejected			0.0250	"	"	"	"	"
cis-1,3-Dichloropropene	U			0.0250	"	"	"	"	"
4-Methyl-2-pentanone	Rejected			0.0250	"	"	"	"	"
trans-1,3-Dichloropropene	U			0.0250	"	"	"	"	"
1,1,2-Trichloroethane	U			0.0250	"	"	"	"	"
Dibromochloromethane	U	J		0.0250	"	"	"	"	"
Toluene	0.618			0.0250	"	"	"	"	"
2-Hexanone	Rejected			0.0250	"	"	"	"	"



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Air Toxics by GC/MS, EPA TO-15 (modified) US EPA Region 5 Chicago Regional Laboratory

1 (1409007-02) Air Sampled: Sep-18-14 10:25 Received: Sep-18-14 14:00

		Flags /		Reporting					
Analyte	Result	Qualifiers	MDL	Limit	Units	Dilution	Batch	Prepared	Analyzed
1,2-Dibromoethane (EDB)	U			0.0250	ppbv	1	B409106	Sep-19-14	Sep-19-14
Tetrachloroethene	U			0.0250	"	"	"	"	"
Chlorobenzene	U			0.0250	"	"	"	"	"
Ethylbenzene	0.0876			0.0250	"	"	"	"	"
m+p-Xylene	0.249			0.100	"	"	"	"	"
Bromoform	U			0.100	"	"	"	"	"
Styrene	U			0.100	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			0.0250	"	"	"	"	"
o-Xylene	0.0911			0.0500	"	"	"	"	"
4-ethyltoluene	0.117	J		0.0250	"	"	"	"	"
1,3,5-Trimethylbenzene	U			0.0500	"	"	"	"	"
1,2,4-Trimethylbenzene	0.160			0.0250	"	"	"	"	"
1,3-Dichlorobenzene	U			0.100	"	"	"	"	"
Benzyl chloride	U			0.0250	"	"	"	"	"
1,4-Dichlorobenzene	U			0.100	"	"	"	"	"
1,2-Dichlorobenzene	U			0.0500	"	"	"	"	"
1,2,4-Trichlorobenzene	0.186			0.100	"	"	"	"	"
Naphthalene	1.15	J, K		0.100	"	"	"	"	"

Surogate	Result		%REC	%REC Limits	Batch	Prepared	Analyzed	
Surrogate: Dichlorodifluoromethane	0.408		78%	90-120	"	"	"	
Surrogate: Dichlorotetrafluoroethane	0.0184		92%	80-120	"	"	"	
Surrogate: Trichlorofluoromethane	0.200		84%	90-120	"	"	"	
Surrogate: 1,1,2-trichloro-1,2,2-trifluoroethane (Fi	0.0346	J	47%	90-120	"	"	"	
Surrogate: 1,1,1-Trichloroethane	5.80E-3	J	116%	80-120	"	"	"	
Surrogate: Carbon tetrachloride	0.0658		77%	80-120	"	"	"	



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Air Toxics by GC/MS, EPA TO-15 (modified) US EPA Region 5 Chicago Regional Laboratory

3 (1409007-03) Air Sampled: Sep-18-14 11:20 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Propene	0.903	K		0.0250	ppbv	1	B409106	Sep-19-14	Sep-19-14
Chloromethane	0.461			0.0250	"	"	"	"	"
Vinyl chloride	U			0.0250	"	"	"	"	"
1,3-butadiene	0.0260	K		0.0250	"	"	"	"	"
Bromomethane	U			0.0250	"	"	"	n .	"
Chloroethane	U			0.0250	"	"	"	n .	"
Acrolein	0.437	K		0.0250	"	"	"	"	"
Isopropyl alcohol	1.09	J		0.0500	"	"	"	"	"
1,1-Dichloroethene	U	J		0.0250	"	"	"	n .	"
Methylene chloride	0.159			0.100	"	"	"	"	"
Carbon disulfide	Rejected			0.0250	"	"	"	"	"
Methyl tert-butyl ether	U			0.0250	"	"	"	"	"
1,1-Dichloroethane	U			0.0250	"	"	"	"	"
Vinyl acetate	U			0.0250	"	"	"	"	"
Hexane	0.366			0.0250	"	"	"	"	"
2-Butanone	0.462			0.0250	"	"	"	"	"
cis-1,2-Dichloroethene	U			0.0250	"	"	"	"	"
trans-1,2-Dichloroethene	U	J		0.0250	"	"	"	"	"
Chloroform	0.0266			0.0250	"	"	"	"	"
1,2-Dichloroethane	U			0.0250	"	"	"	"	"
Cyclohexane	0.109			0.0250	"	"	"	"	"
Benzene	0.200			0.0250	"	"	"	"	"
n-Heptane	0.131	J		0.0250	"	"	"	"	"
1,2-Dichloropropane	U			0.0250	"	"	"	"	"
Trichloroethene	U			0.0250	"	"	"	"	"
Bromodichloromethane	U			0.0250	"	"	"	"	"
1,4-Dioxane	Rejected			0.0250	"	"	"	"	"
cis-1,3-Dichloropropene	U			0.0250	"	"	"	"	"
4-Methyl-2-pentanone	Rejected			0.0250	"	"	"	"	"
trans-1,3-Dichloropropene	U			0.0250	"	"	"	"	"
1,1,2-Trichloroethane	U			0.0250	"	"	"	"	"
Dibromochloromethane	U	J		0.0250	"	"	"	"	"
Toluene	0.433			0.0250	"	"	"	"	"
2-Hexanone	Rejected			0.0250	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			0.0250	"	"	"	"	"
Tetrachloroethene	U			0.0250	"	"	"	"	"



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3 (1409007-03) Air Sampled: Sep-18-14 11:20 Received: Sep-18-14 14:00

		Flags /		Reporting					
Analyte	Result	Qualifiers	MDL	Limit	Units	Dilution	Batch	Prepared	Analyzed
Chlorobenzene	U			0.0250	ppbv	1	B409106	Sep-19-14	Sep-19-14
Ethylbenzene	0.0620			0.0250	"	"	"	"	"
m+p-Xylene	0.177			0.100	"	"	"	"	"
Bromoform	U			0.100	"	"	"	"	"
Styrene	U			0.100	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			0.0250	"	"	"	"	"
o-Xylene	0.0646			0.0500	"	"	"	"	"
4-ethyltoluene	0.0804	J		0.0250	"	"	"	"	"
1,3,5-Trimethylbenzene	U			0.0500	"	"	"	"	"
1,2,4-Trimethylbenzene	0.109			0.0250	"	"	"	"	"
1,3-Dichlorobenzene	U			0.100	"	"	"	"	"
Benzyl chloride	0.0252			0.0250	"	"	"	"	"
1,4-Dichlorobenzene	U			0.100	"	"	"	"	"
1,2-Dichlorobenzene	U			0.0500	"	"	"	"	"
1,2,4-Trichlorobenzene	0.209			0.100	"	"	"	"	"
Naphthalene	1.27	J, K		0.100	"	"	"	"	"

Surogate	Result		%REC	%REC Limits	Batch	Prepared	Analyzed
Surrogate: Dichlorodifluoromethane	0.400		76%	90-120	"	"	"
Surrogate: Dichlorotetrafluoroethane	0.0190		95%	80-120	"	"	"
Surrogate: Trichlorofluoromethane	0.195		82%	90-120	"	"	"
Surrogate: 1,1,2-trichloro-1,2,2-trifluoroethane (Fi	0.0371	J	51%	90-120	"	"	"
Surrogate: 1,1,1-Trichloroethane	5.33E-3	J	107%	80-120	"	"	"
Surrogate: Carbon tetrachloride	0.0674		78%	80-120	"	"	"

3 (1409007-03RE1) Air Sampled: Sep-18-14 11:20 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Ethanol	2.26	J		0.500	ppbv	10	B411044	Sep-22-14	Sep-22-14
Acetone	2.80	J		0.250	"	"	"	"	"

Surogate	Result	%REC	%REC Limits	Batch	Prepared	Analyzed
Surrogate: Dichlorodifluoromethane	0.442	84%	90-120	"	"	"
Surrogate: Trichlorofluoromethane	U	76%	90-120	"	"	"
Surrogate: Carbon tetrachloride	U	77%	80-120	"	"	"



Environmental Protection Agency Region 5

Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605 Phone:(312)353-8370 Fax:(312)886-2591

Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) US EPA Region 5 Chicago Regional Laboratory

2 (1409007-04) Air Sampled: Sep-18-14 12:10 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Propene	1.03	K		0.0250	ppbv	1	B409106	Sep-19-14	Sep-19-14
Chloromethane	0.567			0.0250	"	"	"	"	"
Vinyl chloride	U			0.0250	"	"	"	"	"
1,3-butadiene	0.0392	K		0.0250	"	"	"	"	"
Bromomethane	U			0.0250	"	"	"	"	"
Chloroethane	U			0.0250	"	"	"	"	"
Acrolein	0.722	K		0.0250	"	"	"	"	"
1,1-Dichloroethene	U	J		0.0250	"	"	"	"	"
Carbon disulfide	Rejected			0.0250	"	"	"	"	"
Methyl tert-butyl ether	U			0.0250	"	"	"	"	"
1,1-Dichloroethane	U			0.0250	"	"	"	"	"
Vinyl acetate	U			0.0250	"	"	"	"	"
Hexane	0.512			0.0250	"	"	"	"	"
2-Butanone	0.588			0.0250	"	"	"	"	"
cis-1,2-Dichloroethene	U			0.0250	"	"	"	"	"
trans-1,2-Dichloroethene	U	J		0.0250	"	"	"	"	"
Chloroform	0.0703			0.0250	"	"	"	"	"
1,2-Dichloroethane	U			0.0250	"	"	"	"	"
Cyclohexane	0.156			0.0250	"	"	"	"	"
Benzene	0.300			0.0250	"	"	"	"	"
n-Heptane	0.186	J		0.0250	"	"	"	"	"
1,2-Dichloropropane	U			0.0250	"	"	"	"	"
Trichloroethene	U			0.0250	"	"	"	"	"
Bromodichloromethane	U			0.0250	"	"	"	"	"
1,4-Dioxane	Rejected			0.0250	"	"	"	"	"
cis-1,3-Dichloropropene	U			0.0250	"	"	"	"	"
4-Methyl-2-pentanone	Rejected			0.0250	"	"	"	"	"
trans-1,3-Dichloropropene	U			0.0250	"	"	"	"	"
1,1,2-Trichloroethane	U			0.0250	"	"	"	"	"
Dibromochloromethane	U	J		0.0250	"	"	"	"	"
Toluene	0.679			0.0250	"	"	"	"	"
2-Hexanone	Rejected			0.0250	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			0.0250	"	"	"	"	"
Tetrachloroethene	U			0.0250	"	"	"	"	"
Chlorobenzene	U			0.0250	"	"	"	"	"
Ethylbenzene	0.149			0.0250	"	"	"	"	"



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Air Division, US EPA Region 5 Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) US EPA Region 5 Chicago Regional Laboratory

2 (1409007-04) Air Sampled: Sep-18-14 12:10 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
m+p-Xylene	0.473			0.100	ppbv	1	B409106	Sep-19-14	Sep-19-14
Bromoform	U			0.100	"	"	"	"	"
Styrene	U			0.100	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			0.0250	"	"	"	"	"
o-Xylene	0.162			0.0500	"	"	"	"	"
4-ethyltoluene	0.145	J		0.0250	"	"	"	"	"
1,3,5-Trimethylbenzene	U			0.0500	"	"	"	"	"
1,2,4-Trimethylbenzene	0.190			0.0250	"	"	"	"	"
1,3-Dichlorobenzene	U			0.100	"	"	"	"	"
Benzyl chloride	0.0704			0.0250	"	"	"	"	"
1,4-Dichlorobenzene	U			0.100	"	"	"	"	"
1,2-Dichlorobenzene	U			0.0500	"	"	"	"	"
1,2,4-Trichlorobenzene	0.190			0.100	"	"	"	"	"
Naphthalene	0.973	J, K		0.100	"	"	"	"	"

Surogate	Result		%REC	%REC Limits	Batch	Prepared	Analyzed
Surrogate: Dichlorodifluoromethane	0.425		81%	90-120	"	"	"
Surrogate: Dichlorotetrafluoroethane	0.0209		104%	80-120	"	"	"
Surrogate: Trichlorofluoromethane	0.295		123%	90-120	"	"	"
Surrogate: 1,1,2-trichloro-1,2,2-trifluoroethane (Fi	0.0471	J	65%	90-120	"	"	"
Surrogate: 1,1,1-Trichloroethane	6.28E-3	J	126%	80-120	"	"	"
Surrogate: Carbon tetrachloride	0.123		143%	80-120	"	"	"

2 (1409007-04RE1) Air Sampled: Sep-18-14 12:10 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Isopropyl alcohol	1.02	J		0.0500	ppbv	1	B411044	Sep-23-14	Sep-23-14
Methylene chloride	0.164	K		0.100	"	"	"	"	"

Surogate	Result	%REC	%REC Limits	Batch	Prepared	Analyzed
Surrogate: Dichlorodifluoromethane	0.397	76%	90-120	"	"	<i>"</i>
Surrogate: Trichlorofluoromethane	0.200	84%	90-120	"	"	"
Surrogate: Carbon tetrachloride	0.0694	81%	80-120	"	"	"



Chicago Regional Laboratory

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Air Division, US EPA Region 5 Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) US EPA Region 5 Chicago Regional Laboratory

2 (1409007-04RE2) Air Sampled: Sep-18-14 12:10 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Ethanol	2.34	J		0.333	ppbv	6.667	B411044	Sep-22-14	Sep-22-14
Acetone	2.42	J, K		0.167	"	"	"	"	"

Surogate	Result	%REC	%REC Limits	Batch	Prepared	Analyzed
Surrogate: Dichlorodifluoromethane	0.483	92%	90-120	"	"	"
Surrogate: Trichlorofluoromethane	U	81%	90-120	"	"	"
Surrogate: Carbon tetrachloride	U	67%	80-120	"	"	"

4B (1409007-05) Air Sampled: Sep-18-14 00:00 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Propene	0.927	K		0.0250	ppbv	1	B409106	Sep-19-14	Sep-19-14
Chloromethane	0.506			0.0250	"	"	"	"	"
Vinyl chloride	U			0.0250	"	"	"	"	"
1,3-butadiene	U	K		0.0250	"	"	"	"	"
Bromomethane	U			0.0250	"	"	"	"	"
Chloroethane	U			0.0250	"	"	"	"	"
Acrolein	0.531	K		0.0250	"	"	"	"	"
Isopropyl alcohol	1.14	J		0.0500	"	"	"	"	"
1,1-Dichloroethene	U	J		0.0250	"	"	"	"	"
Methylene chloride	0.163	K		0.100	"	"	"	"	"
Carbon disulfide	Rejected			0.0250	"	"	"	"	"
Methyl tert-butyl ether	U			0.0250	"	"	"	"	"
1,1-Dichloroethane	U			0.0250	"	"	"	"	"
Vinyl acetate	U			0.0250	"	"	"	"	"
Hexane	0.391			0.0250	"	"	"	"	"
2-Butanone	0.410			0.0250	"	"	"	"	"
cis-1,2-Dichloroethene	U			0.0250	"	"	"	"	"
trans-1,2-Dichloroethene	U	J		0.0250	"	"	"	"	"
Chloroform	0.0255			0.0250	"	"	"	"	"
1,2-Dichloroethane	U			0.0250	"	"	"	"	"
Cyclohexane	0.126			0.0250	"	"	"	"	"
Benzene	0.230			0.0250	"	"	"	"	"
n-Heptane	0.148	J		0.0250	"	"	"	"	"
1,2-Dichloropropane	U			0.0250	"	"	"	"	"



Surrogate: 1,1,1-Trichloroethane

Surrogate: Carbon tetrachloride

3.45E-3

0.0624

Air Division, US EPA Region 5

Environmental Protection Agency Region 5

Chicago Regional Laboratory

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Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) US EPA Region 5 Chicago Regional Laboratory

4B (1409007-05) Air Sampled: Sep-18-14 00:00 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Trichloroethene	U			0.0250	ppbv	1	B409106	Sep-19-14	Sep-19-14
Bromodichloromethane	U			0.0250	"	"	"	"	"
1,4-Dioxane	Rejected			0.0250	"	"	"	"	"
cis-1,3-Dichloropropene	U			0.0250	"	"	"	"	"
4-Methyl-2-pentanone	Rejected			0.0250	"	"	"	"	"
trans-1,3-Dichloropropene	U			0.0250	"	"	"	"	"
1,1,2-Trichloroethane	U			0.0250	"	"	"	"	"
Dibromochloromethane	U	J		0.0250	"	"	"	"	"
Toluene	0.462			0.0250	"	"	"	"	"
2-Hexanone	Rejected			0.0250	"	"	"	"	"
1,2-Dibromoethane (EDB)	U			0.0250	"	"	"	"	"
Tetrachloroethene	U			0.0250	"	"	"	"	"
Chlorobenzene	U			0.0250	"	"	"	"	"
Ethylbenzene	0.0689			0.0250	"	"	"	"	"
m+p-Xylene	0.185			0.100	"	"	"	"	"
Bromoform	U			0.100	"	"	"	"	"
Styrene	U			0.100	"	"	"	"	"
1,1,2,2-Tetrachloroethane	U			0.0250	"	"	"	"	"
o-Xylene	0.0702			0.0500	"	"	"	"	"
4-ethyltoluene	0.0755	J		0.0250	"	"	"	"	"
1,3,5-Trimethylbenzene	U			0.0500	"	"	"	"	"
1,2,4-Trimethylbenzene	0.0898			0.0250	"	"	"	"	"
1,3-Dichlorobenzene	U			0.100	"	"	"	"	"
Benzyl chloride	U			0.0250	"	"	"	"	"
1,4-Dichlorobenzene	U			0.100	"	"	"	n .	"
1,2-Dichlorobenzene	U			0.0500	"	"	"	"	"
1,2,4-Trichlorobenzene	0.177			0.100	"	"	"	n .	"
Naphthalene	1.00	J, K		0.100	"	"	"	"	"
Surogate	Result			%REC		%REC Limits	Batch	Prepared	Analyzed
Surrogate: Dichlorodifluoromethane	0.426			81%		90-120	"	"	"
Surrogate: Dichlorotetrafluoroethane	0.0203			102%		80-120	"	"	"
Surrogate: Trichlorofluoromethane	0.202			85%		90-120	"	"	"
Surrogate: 1,1,2-trichloro-1,2,2-trifluoroethane (Fi	0.0367	J		50%		90-120	"	"	"

69%

73%

80-120

80-120



Environmental Protection Agency Region 5 Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605

Project: BP Whiting Refinery

Phone:(312)353-8370 Fax:(312)886-2591

77 West Jackson Boulevard Project Number: [none] Reported: Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) **US EPA Region 5 Chicago Regional Laboratory**

4B (1409007-05RE2) Air Sampled: Sep-18-14 00:00 Received: Sep-18-14 14:00

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Ethanol	3.34	J		0.333	ppbv	6.667	B411044	Sep-22-14	Sep-22-14
Acetone	2.39	K, J		0.167	"	"	"	"	"

Surogate	Result	%REC	%REC Limits	Batch	Prepared	Analyzed
Surrogate: Dichlorodifluoromethane	0.448	85%	90-120	"	"	"
Surrogate: Trichlorofluoromethane	U	79%	90-120	"	"	"
Surrogate: Carbon tetrachloride	U	65%	80-120	"	"	"

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Air Division, US EPA Region 5 Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B409106 - ColdTrap Dehydration

Propose	Blank (B409106-BLK1)	Prepared & Analyzed: Sep-19-14										
Propene U 0.0250 pbv 1.1.					Reporting		Spike	Source		%REC		RPD
Chloromethane U 0.0250 " Vinyl chloride U 0.0250 " Bromomethane U 0.0250 " Chloroethane U 0.0250 " Ethanol U 0.0250 " Acrolein 0.028 0.0500 " Ropropyl alcohol 0.108 J 0.0500 " Actone 0.106 J 0.0250 " Actone 0.106 J 0.0250 " Li-Dichloroethane U 0.0250 " Methyl tert-butyl ether U 0.0250 " Uilyl acetate U 0.0250 " Vinyl acetate U 0.0250 " Hexane U 0.0250 " 2-Butanone U 0.0250 " 5-Indichorethane U 0.0250 " Chloroforn U 0.0250 " Cyclohexane U 0.0250	Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit
Strong claims U	Propene	U			0.0250	ppbv						
National	Chloromethane	U			0.0250	"						
Chlorocthane	Vinyl chloride	U			0.0250	"						
Chloroethane U J.0.0250 " Ethanol U.0.258 0.0580 " Acrolein 0.0258 0.0580 " Sopropyl alcohol 0.105 J.0.0500 " Acetone 0.106 J.0.0250 " Methylenchloride U.0.0250 " Methyl terr-butyl ether U.0.0250 " J.1-Dichloroethane U.0.0250 " Winyl acetate U.0.0250 " Hexane U.0.0250 " 2-Butanone U.0.0250 " cis-1,2-Dichloroethene U.0.0250 " Urans-1,2-Dichloroethene U.0.0250 " Chloroform U.0.0250 " Cyclohexane U.0.0250 " Gyclohexane U.0.0250 " Benzee U.0.0250 " 1-L-Dichloropropane U.0.0250 " Trickloroethane U.0.0250 " Bromodichloromethane U.0.0250 " <th< th=""><th>1,3-butadiene</th><th>U</th><th></th><th></th><th>0.0250</th><th>"</th><th></th><th></th><th></th><th></th><th></th><th></th></th<>	1,3-butadiene	U			0.0250	"						
Ethanol U J 0.0500 " Acrolein 0.0258 0.0250 " Isopropyl alcohol 0.106 J 0.0500 " Acteone 0.106 J 0.0250 " Herbylen chloride U J 0.0250 " Methyl tert-butyl ether U 0.0250 " Heydle tert-butyl ether U 0.0250 " Hill-Dichloroethane U 0.0250 " Vinyl acetate U 0.0250 " Hexane U 0.0250 " 2-Butanone U 0.0250 " 1-Li-Dichloroethene U 0.0250 "	Bromomethane	U			0.0250	"						
Acrolein	Chloroethane	U			0.0250	"						
Sopropyl alcohol 0.105 3 0.050 5 7 0.0250 5 7 0.0250 5 7 0.0250 5 7 0.0250 5 7 0.0250 7 0.025	Ethanol	U	J		0.0500	"						
Actone 0.106 J 0.0250 " Methylene chloride U J 0.0250 " Carbon disulfide Rejected 0.0250 " Methyl tert-butyl ether U 0.0250 " I,1-Dichloroethane U 0.0250 " Vinyl acetate U 0.0250 " Hexane U 0.0250 " 2-Butanone U 0.0250 " cis-1,2-Dichloroethene U 0.0250 " trans-1,2-Dichloroethene U 0.0250 " Chloroform U 0.0250 " Cyclohexane U 0.0250 " Benzene U 0.0250 " n-Heptane U 0.0250 " 1,2-Dichloropropane U 0.0250 " Bromodichloromethane U 0.0250 " Bromodichloropropene U 0.0250 " 4-Methyl-2-pentanone R	Acrolein	0.0258			0.0250	"						
1,1-Dichloroethene	Isopropyl alcohol	0.105	J		0.0500	"						
Methylene chloride Carbon disulfide Rejected Methylene chloride U 0.0250 " 1,1-Dichloroethane U 0.0250 " Vinyl acetate U 0.0250 " Hexane U 0.0250 " 1,2-Dichloroethene U 0.0250 " 0.0250	Acetone	0.106	J		0.0250	"						
Carbon disulfide Rejected 0.0250 " Methyl tert-butyl ether U 0.0250 " 1,1-Dichloroethane U 0.0250 " Vinyl acetate U 0.0250 " Hexane U 0.0250 " 2-Butanone U 0.0250 " cis-1,2-Dichloroethene U 0.0250 " Chloroform U 0.0250 " 1,2-Dichloroethane U 0.0250 " Cyclohexane U 0.0250 " Benzene U 0.0250 " 1,2-Dichloropropane U 0.0250 " 1,2-Dichloropropane U 0.0250 " Bromodichloromethane U 0.0250 " Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " 4-Methyl-2-pentanone U <th< th=""><th>1,1-Dichloroethene</th><th>U</th><th>J</th><th></th><th>0.0250</th><th>"</th><th></th><th></th><th></th><th></th><th></th><th></th></th<>	1,1-Dichloroethene	U	J		0.0250	"						
Methyl tert-butyl ether U 0.0250 " 1,1-Dichloroethane U 0.0250 " Vinyl acetate U 0.0250 " Hexane U 0.0250 " 2-Butanone U 0.0250 " cis-1,2-Dichloroethene U 0.0250 " trans-1,2-Dichloroethene U J 0.0250 " Chloroform U 0.0250 " 1,2-Dichloroethane U 0.0250 " Cyclohexane U 0.0250 " Benzene U 0.0250 " 1-2-Dichloropropane U 0.0250 " Trichloroethene U 0.0250 " Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " 4-Methyl-2-pentanone U 0.0250 "	Methylene chloride	U			0.100	"						
1,1-Dichloroethane	Carbon disulfide	Rejected			0.0250	"						
Vinyl acetate U 0.0250 " Hexane U 0.0250 " 2-Butanone U 0.0250 " cis-1,2-Dichloroethene U 0.0250 " Chloroform U 0.0250 " 1,2-Dichloroethane U 0.0250 " Cyclohexane U 0.0250 " Benzene U 0.0250 " 1,2-Dichloropropane U 0.0250 " 1,2-Dichloropropane U 0.0250 " Trichloroethene U 0.0250 " Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 "	Methyl tert-butyl ether	U			0.0250	"						
	1,1-Dichloroethane	U			0.0250	"						
Description	Vinyl acetate	U			0.0250	"						
cis-1,2-Dichloroethene U 0.0250 " trans-1,2-Dichloroethene U J 0.0250 " Chloroform U 0.0250 " 1,2-Dichloroethane U 0.0250 " Cyclohexane U 0.0250 " Benzene U 0.0250 " n-Heptane U J 0.0250 " 1,2-Dichloropropane U 0.0250 " Trichloroethene U 0.0250 " Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " cis-1,3-Dichloropropene U 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	Hexane	U			0.0250	"						
trans-1,2-Dichloroethene U J 0.0250 " Chloroform U 0.0250 " 1,2-Dichloroethane U 0.0250 " Cyclohexane U 0.0250 " Benzene U J 0.0250 " n-Heptane U J 0.0250 " 1,2-Dichloropropane U 0.0250 " Trichloroethene U 0.0250 " Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	2-Butanone	U			0.0250	"						
Chloroform U 0.0250 " 1,2-Dichloroethane U 0.0250 " Cyclohexane U 0.0250 " Benzene U 0.0250 " n-Heptane U 0.0250 " 1,2-Dichloropropane U 0.0250 " Trichloroethene U 0.0250 " Trichloroethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " 1,4-Methyl-2-pentanone Rejected U 0.0250 " 1,4-Methyl-2-pentanone U 0.0250 " 1,4-Methyl-	cis-1,2-Dichloroethene	U			0.0250	"						
1,2-Dichloroethane U 0.0250 " Cyclohexane U 0.0250 " Benzene U 0.0250 " n-Heptane U J 0.0250 " 1,2-Dichloropropane U 0.0250 " Trichloroethene U 0.0250 " Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " cis-1,3-Dichloropropene U 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	trans-1,2-Dichloroethene	U	J		0.0250	"						
Cyclohexane U 0.0250 " Benzene U 0.0250 " n-Heptane U J 0.0250 " 1,2-Dichloropropane U 0.0250 " Trichloroethene U 0.0250 " Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " cis-1,3-Dichloropropene U 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	Chloroform	U			0.0250	"						
Benzene	1,2-Dichloroethane	U			0.0250	"						
n-Heptane U J 0.0250 " 1,2-Dichloropropane U 0.0250 " Trichloroethene U 0.0250 " Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " cis-1,3-Dichloropropene U 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	Cyclohexane	U			0.0250	"						
1,2-Dichloropropane U 0.0250 " Trichloroethene U 0.0250 " Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " cis-1,3-Dichloropropene U 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	Benzene	U			0.0250	"						
Trichloroethene U 0.0250 " Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " cis-1,3-Dichloropropene U 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	n-Heptane	U	J		0.0250	"						
Bromodichloromethane U 0.0250 " 1,4-Dioxane Rejected 0.0250 " cis-1,3-Dichloropropene U 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	1,2-Dichloropropane	U			0.0250	"						
1,4-Dioxane Rejected 0.0250 " cis-1,3-Dichloropropene U 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	Trichloroethene	U			0.0250	"						
cis-1,3-Dichloropropene U 0.0250 " 4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	Bromodichloromethane	U			0.0250	"						
4-Methyl-2-pentanone Rejected 0.0250 " trans-1,3-Dichloropropene U 0.0250 "	1,4-Dioxane	Rejected			0.0250	"						
trans-1,3-Dichloropropene U 0.0250 "	cis-1,3-Dichloropropene	U			0.0250	"						
	4-Methyl-2-pentanone	Rejected			0.0250	"						
1,1,2-Trichloroethane U 0.0250 "	trans-1,3-Dichloropropene	U			0.0250	"						
	1,1,2-Trichloroethane	U			0.0250	"						



Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605 Phone:(312)353-8370 Fax:(312)886-2591

Air Division, US EPA Region 5 Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B409106 - ColdTrap Dehydration

Blank (B409106-BLK1)	Prepared & Analyzed: Sep-19-14											
		Flags /		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	
Dibromochloromethane	U	J		0.0250	ppbv							
Toluene	U			0.0250	"							
2-Hexanone	Rejected			0.0250	"							
1,2-Dibromoethane (EDB)	U			0.0250	"							
Tetrachloroethene	U			0.0250	"							
Chlorobenzene	U			0.0250	"							
Ethylbenzene	U			0.0250	"							
m+p-Xylene	U			0.100	"							
Bromoform	U			0.100	"							
Styrene	U			0.100	"							
1,1,2,2-Tetrachloroethane	U			0.0250	"							
o-Xylene	U			0.0500	"							
4-ethyltoluene	U	J		0.0250	"							
1,3,5-Trimethylbenzene	U			0.0500	"							
1,2,4-Trimethylbenzene	U			0.0250	"							
1,3-Dichlorobenzene	U			0.100	"							
Benzyl chloride	U			0.0250	"							
1,4-Dichlorobenzene	U			0.100	"							
1,2-Dichlorobenzene	U			0.0500	"							
1,2,4-Trichlorobenzene	U			0.100	"							
Naphthalene	U	J		0.100	"							
Surrogate: Dichlorodifluoromethane	U				"	0.5250		%	90-120			
Surrogate: Dichlorotetrafluoroethane	3.85E-3				"	2.000E-2		19%	80-120			
Surrogate: Trichlorofluoromethane	U				"	0.2390		%	90-120			
Surrogate:	U	J			"	7.300E-2		%	90-120			
1,1,2-trichloro-1,2,2-trifluoroethane (Freon												
Surrogate: 1,1,1-Trichloroethane	3.05E-3	J			"	5.000E-3		61%	80-120			
Surrogate: Carbon tetrachloride	U				"	8.600E-2		%	80-120			



Chicago Regional Laboratory

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Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B409106 - ColdTrap Dehydration

Propose	LCS (B409106-BS1)		Prepared & Analyzed: Sep-19-14										
Propose			•		Reporting		Spike	Source		%REC		RPD	
Chloromethane 0.507 0.0250 0.5000 101% 77.6-116 Vinyl chloride 0.644 0.0250 0.5000 101% 77.4-117 Ja-butdiene 0.643 0.0250 0.5000 93% 77.7-115 Bromomethane 0.644 0.0250 0.5000 93% 78.2-117 Chlorocthane 0.665 0.0250 0.5000 93% 78.2-117 Ethanol 0.672 J 0.0500 0.5000 93% 679-130 Isopropyl alcohol 0.329 J 0.0500 0.5000 98% 679-130 Actione 0.442 J 0.0250 0.5000 88% 61-137 Hethylene chloride 0.431 J 0.0250 0.5000 88% 61-137 Methylene chloride 0.521 0.100 0.5000 99% 73-118 Methylene chloride 0.431 0.0250 0.5000 89% 79-118 Li-Dichlorothane 0.431 0.0250 0.5000 <t< th=""><th>Analyte</th><th>Result</th><th>Qualifiers</th><th>MDL</th><th>Limit</th><th>Units</th><th>Level</th><th>Result</th><th>%REC</th><th>Limits</th><th>RPD</th><th>Limit</th></t<>	Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	
Name	Propene	0.511			0.0250	ppbv	0.5000		102%	76-112			
1,-butaliene	Chloromethane	0.507			0.0250	"	0.5000		101%	77.6-116			
Bromomethane	Vinyl chloride	0.504			0.0250	"	0.5000		101%	77.4-117			
Chlorocthane 0.465 Chlorocthane 0.4672 J 0.0500 - 0.5000 39% 79.5-114 Chlorocthane 0.470 0.250 - 0.5000 39% 79.5-114 Chlorocthane 0.480 0.250 - 0.5000 39% 79.5-114 Chlorocthane 0.412 J 0.0500 - 0.5000 39% 609-137 L1-Dichlorocthene 0.442 J 0.0500 - 0.5000 38% 60-137 L1-Dichlorocthene 0.445 J 0.0250 - 0.5000 38% 60-137 L1-Dichlorocthene 0.445 - 0.450 - 0.5000 38% 77.3-118 Methylere-thoride 0.521 0.0250 - 0.5000 38% 77.3-118 Methylere-thutyl ether 0.409 - 0.0250 - 0.5000 38% 79.5-117 Chlorocthane 0.431 0.0250 - 0.0500 38% 79.5-117 Vinyl acetate 0.437 0.0250 - 0.5000 38% 79.5-117 Vinyl acetate 0.441 0.0250 - 0.5000 38% 79.5-118 Cyclohexane 0.441 0.0250 - 0.5000 0.500	1,3-butadiene	0.463			0.0250	"	0.5000		93%	77.7-115			
Ethanlo 0.672 J 0.0300 ° 0.5000 134% 50-124 Acrolein 0.490 0.0250 ° 0.5000 .98% 67.9-130 Isopropi alcohol 0.329 J 0.0500 ° 0.5000 .88% 67.9-130 Acceone 0.442 J 0.0250 ° 0.5000 .88% 60-137 I.1-Dichloroethene 0.445 J 0.0250 ° 0.5000 .99% 73118 Methylene chloride 0.521 0.0250 ° 0.5000 .99% 73118 Methylete-butje ther 0.499 0.0250 ° 0.5000 .99% 73118 Heyane 0.431 0.0250 ° 0.5000 .98% 79118 Heyane 0.431 0.0250 ° 0.5000 .88% 79118 Leyane 0.431 0.0250 ° 0.5000 .28% 70121 Eblation or 0.460 0.0250 ° <th< th=""><th>Bromomethane</th><th>0.504</th><th></th><th></th><th>0.0250</th><th>"</th><th>0.5000</th><th></th><th>101%</th><th>75.2-117</th><th></th><th></th></th<>	Bromomethane	0.504			0.0250	"	0.5000		101%	75.2-117			
Acrolein 0.490 0.0250 0.0500 98% 67.9-130 Isopropyl alcohol 0.329 J 0.0500 0.5000 66% 81-106 Acctone 0.442 J 0.0250 0.5000 88% 60-137 Hetholeroethene 0.445 J 0.0250 0.5000 89% 77.3-118 Methylne chloride 0.521 1.00 0.0250 0.5000 89% 73.118 Methylne chloride 0.621 0.0250 0.5000 90% 53.140 Methylne chloride 0.490 0.0250 0.5000 88% 79.118 I,1-Dichloroethane 0.431 0.0250 0.5000 88% 79.118 Hexane 0.437 0.0250 0.5000 88% 79.117 2-Butanone 0.363 0.0250 0.5000 37% 85.106 6is-1,2-Dichloroethane 0.460 0.0250 0.5000 37% 78.115 Chloroform 0.491 0.0250 0.5000 98%<	Chloroethane	0.465			0.0250	"	0.5000		93%	79.8-114			
Sopropyl alcohol 0.329 J 0.0500 0.5000 66% 81-106 Acetone 0.442 J 0.0250 0.5000 88% 60-137 I.1-Dichloroethene 0.445 J 0.0250 0.5000 89% 77.3-118 Methylne chloride 0.521 0.1000 0.5000 0.5000 90% 53-140 Methyl tert-butyl ether 0.409 0.0250 0.5000 0.5000 86% 78.8-117 Methyl tert-butyl ether 0.431 0.0250 0.5000 86% 79.8-117 Winyl acetate 0.373 0.0250 0.5000 86% 79.8-117 Winyl acetate 0.437 0.0250 0.5000 87% 75.118 Hexane 0.437 0.0250 0.5000 75% 85-106 Hexane 0.437 0.0250 0.5000 75% 85-106 Hexane 0.460 0.5000 0.5000 75% 82-108 Cis-1,2-Dichloroethene 0.460 0.0250 0.5000 0.5000 75% 82-108 Cis-1,2-Dichloroethene 0.461 0.0250 0.5000 0.5000 0.75% 82-108 Cis-1,2-Dichloroethene 0.491 0.0250 0.5000 0.5000 0.75% 82-108 Cis-1,2-Dichloroethene 0.491 0.0250 0.5000 0.5000 0.5000 0.5000 Cis-1,2-Dichloroethene 0.491 0.0250 0.5000 0.5000 0.5000 0.5000 Cis-1,2-Dichloroethene 0.491 0.0250 0.5000 0.5000 0.5000 0.5000 0.5000 Cis-1,2-Dichloroethene 0.491 0.0250 0.5000 0.5000 0.5000 0.5000 0.5000 0.5000 Cis-1,2-Dichloroethene 0.491 0.0250 0.5000 0.5000 0.5000 0.5000 0.5000 0.5000 0.5000 Cis-1,2-Dichlorophone 0.494 0.0250 0.5000 0	Ethanol	0.672	J		0.0500	"	0.5000		134%	50-124			
Actetone 0.442 J 0.0250 " 0.5000 88% 60-137 1,1-Dichloroethene 0.445 J 0.0250 " 0.5000 88% 60-137 1,1-Dichloroethene 0.521 0.100 " 0.5000 104% 78.7-117 Carbon disulfide Rejected 0.0250 " 0.5000 90% 53-140 Methyl terr-butyl ether 0.409 0.0250 " 0.5000 82% 79.1-118 1,1-Dichloroethane 0.431 0.0250 " 0.5000 88% 79.8-117 Vinyl acetate 0.373 0.0250 " 0.5000 88% 79.8-117 Vinyl acetate 0.373 0.0250 " 0.5000 87% 75.8-116 Hexane 0.437 0.0250 " 0.5000 77% 85-106 Hexane 0.460 0.0250 " 0.5000 77% 82-108 cis-1,2-Dichloroethene 0.460 0.0250 " 0.5000 92% 78.1-115 trans-1,2-Dichloroethane 0.501 J 0.0250 " 0.5000 98% 79.6-115 1,2-Dichloroethane 0.531 0.0250 " 0.5000 98% 79.6-115 1,2-Dichloroethane 0.441 0.0250 " 0.5000 98% 79.6-115 1,2-Dichloroethane 0.441 0.0250 " 0.5000 98% 79.6-115 Therman dichloropropane 0.532 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.445 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.445 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.446 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.447 J 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.446 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.447 J 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.446 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.447 J 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.446 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.447 J 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.447 J 0.0250 " 0.5000 98% 75.7-118 Trichloroethane 0.445 0.0250 " 0.5000 98% 75.8-117 Trichloroethane 0.446 0.0250 " 0.5000 98% 75.8-117 Trichloroethane 0.446 0.0250 " 0.5000 98% 75.5-115 Trichloroethane 0.447 0.0250 " 0.5000 98% 75.5-115 Trichloroethane 0.446 0.0250 " 0.5000 98% 75.5-115 Trichloroethane 0.447 0.0250 " 0.5000 98% 75.5-115 Trichloroethane 0.448 0.0250 " 0.5000 98% 75.5-115 Trichloroethane 0.448 0.0250 " 0.5000 98% 75.5-115 Trichloroethane 0.448 0.0250 " 0.5000 98% 75.5-115	Acrolein	0.490			0.0250	"	0.5000		98%	67.9-130			
1.1-Dichloroethene	Isopropyl alcohol	0.329	J		0.0500	"	0.5000		66%	81-106			
Methylene chloride 0.521 0.100 0.5000 104% 78.7117 Carbon disulfide Rejected 0.0250 0.5000 90% 53-140 Methyle terl-butyl ether 0.409 0.0250 0.5000 82% 79.1-118 1,1-Dichlorocthane 0.431 0.0250 0.5000 86% 79.8-117 Vinyl acetate 0.373 0.0250 0.5000 87% 70.1-121 Lexane 0.437 0.0250 0.5000 87% 70.1-121 2-Butanone 0.363 0.0250 0.5000 87% 70.1-121 2-Butanone 0.460 0.0250 0.5000 87% 70.1-121 2-Butanone 0.460 0.0250 0.0250 0.5000 92% 78.1-115 trans-1,2-Dichlorocthane 0.501 J 0.0250 0.5000 98% 79.6-115 1,2-Dichlorocthane 0.531 0.0250 0.0250 0.5000 98% 79.5-115 Cyclohexane 0.441 0.0250 0.0250 </th <th>Acetone</th> <th>0.442</th> <th>J</th> <th></th> <th>0.0250</th> <th>"</th> <th>0.5000</th> <th></th> <th>88%</th> <th>60-137</th> <th></th> <th></th>	Acetone	0.442	J		0.0250	"	0.5000		88%	60-137			
Carbon disulfide Rejected 0.0250 " 0.5000 95% 53-140 Methyl tert-butyl ether 0.409 0.0250 " 0.5000 82% 79.1-118 1,1-Dichloroethane 0.431 0.0250 " 0.5000 86% 79.8-117 Vinyl acetate 0.373 0.0250 " 0.5000 87% 70.1-121 2-Butanone 0.437 0.0250 " 0.5000 87% 70.1-121 2-Butanone 0.460 0.0250 " 0.5000 73% 82-108 cis-1,2-Dichloroethane 0.501 J 0.0250 " 0.5000 92% 78.1-115 Ly2-Dichloroethane 0.501 J 0.0250 " 0.5000 98% 796-115 1,2-Dichloroethane 0.531 0.0250 " 0.5000 98% 796-115 Cyclohexane 0.441 0.0250 " 0.5000 98% 79.5-118 Benzene 0.494 0.0250 " <	1,1-Dichloroethene	0.445	J		0.0250	"	0.5000		89%	77.3-118			
Methyl tert-butyl ether 0.409 0.0250 0.5000 82% 79.1-118 1,1-Dichloroethane 0.431 0.0250 0.5000 86% 79.8-117 Vinyl acetate 0.373 0.0250 0.5000 87% 85-106 Hexane 0.437 0.0250 0.5000 87% 70.1-121 2-Butanone 0.363 0.0250 0.5000 73% 82-108 cis-1,2-Dichloroethene 0.460 0.0250 0.5000 92% 78.1-115 Cisn-1,2-Dichloroethene 0.460 0.0250 0.5000 98% 79.6-115 Lip-Dichloroethane 0.501 J 0.0250 0.5000 98% 79.6-115 Lip-Dichloroethane 0.531 0.0250 0.0250 0.5000 98% 79.6-115 Cyclohexane 0.441 0.0250 0.0250 0.5000 88% 75.7-118 Benzene 0.494 0.0250 0.0500 89% 66.9-152 1,2-Dichloropropane 0.445 0.0250 0	Methylene chloride	0.521			0.100	"	0.5000		104%	78.7-117			
1.1-Dichloroethane 0.431 0.0250 0.5000 86% 79.8-117	Carbon disulfide	Rejected			0.0250	"	0.5000		90%	53-140			
No.	Methyl tert-butyl ether	0.409			0.0250	"	0.5000		82%	79.1-118			
Hexane 0.437 0.0250 " 0.5000 87% 70.1-121 2-Butanone 0.363 0.0250 " 0.5000 73% 82-108 cis-1,2-Dichloroethene 0.460 0.0250 " 0.5000 92% 78.1-115 trans-1,2-Dichloroethene 0.501	1,1-Dichloroethane	0.431			0.0250	"	0.5000		86%	79.8-117			
2-Butanone 0.363 0.0250 " 0.5000 73% 82-108 cis-1,2-Dichloroethene 0.460 0.0250 " 0.5000 92% 78.1-115 trans-1,2-Dichloroethene 0.501 J 0.0250 " 0.5000 92% 78.1-115 trans-1,2-Dichloroethene 0.501 J 0.0250 " 0.5000 98% 79.6-115 1,2-Dichloroethane 0.531 0.0250 " 0.5000 98% 79.6-115 1,2-Dichloroethane 0.531 0.0250 " 0.5000 88% 72.5-119	Vinyl acetate	0.373			0.0250	"	0.5000		75%	85-106			
cis-1,2-Dichloroethene 0.460 0.0250 " 0.5000 92% 78.1-115 trans-1,2-Dichloroethene 0.501 J 0.0250 " 0.5000 100% 0-200 Chloroform 0.491 0.0250 " 0.5000 98% 79.6-115 1,2-Dichloroethane 0.531 0.0250 " 0.5000 88% 72.5-119 Benzene 0.441 0.0250 " 0.5000 88% 72.5-119 Benzene 0.447 J 0.0250 " 0.5000 89% 66.9-152 1,2-Dichloropropane 0.532 0.0250 " 0.5000 89% 66.9-152 1,2-Dichloropropane 0.445 0.0250 " 0.5000 89% 70.1-119 Bromodichloromethane 0.445 0.0250 " 0.5000 89% 70.1-119 1,4-Dioxane Rejected 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 <th< th=""><th>Hexane</th><th>0.437</th><th></th><th></th><th>0.0250</th><th>"</th><th>0.5000</th><th></th><th>87%</th><th>70.1-121</th><th></th><th></th></th<>	Hexane	0.437			0.0250	"	0.5000		87%	70.1-121			
trans-1,2-Dichloroethene 0.501 J 0.0250 " 0.5000 100% 0-200 Chloroform 0.491 0.0250 " 0.5000 98% 79.6-115 1,2-Dichloroethane 0.531 0.0250 " 0.5000 88% 79.8-115 Cyclohexane 0.441 0.0250 " 0.5000 88% 72.5-119 Benzene 0.494 0.0250 " 0.5000 89% 66.9-152 1,2-Dichloropropane 0.447 J 0.0250 " 0.5000 89% 66.9-152 1,2-Dichloropropane 0.532 0.0250 " 0.5000 89% 76.8-118 Trichloroethene 0.445 0.0250 " 0.5000 89% 70.1-119 Bromodichloromethane 0.473 0.0250 " 0.5000 81% 54.7-150 cis-1,3-Dichloropropene 0.436 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone 0.463 0.0250	2-Butanone	0.363			0.0250	"	0.5000		73%	82-108			
Chloroform 0.491 0.0250 " 0.5000 98% 79.6-115 1,2-Dichloroethane 0.531 0.0250 " 0.5000 106% 79.8-115 Cyclohexane 0.441 0.0250 " 0.5000 88% 72.5-119 Benzene 0.494 0.0250 " 0.5000 89% 75.7-118 n-Heptane 0.447 J 0.0250 " 0.5000 89% 66.9-152 1,2-Dichloropropane 0.532 0.0250 " 0.5000 89% 70.1-119 Bromodichloromethane 0.445 0.0250 " 0.5000 89% 70.1-119 Horosane Rejected 0.0250 " 0.5000 87% 75.8-117 1,4-Dioxane Rejected 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 " 0.5000 87% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 " 0.5000	cis-1,2-Dichloroethene	0.460			0.0250	"	0.5000		92%	78.1-115			
1,2-Dichloroethane 0.531 0.0250 " 0.5000 106% 79.8-115 Cyclohexane 0.441 0.0250 " 0.5000 88% 72.5-119 Benzene 0.494 0.0250 " 0.5000 99% 75.7-118 n-Heptane 0.447 J 0.0250 " 0.5000 89% 66.9-152 1,2-Dichloropropane 0.532 0.0250 " 0.5000 89% 76.8-118 Trichloroethene 0.445 0.0250 " 0.5000 89% 70.1-119 Bromodichloromethane 0.473 0.0250 " 0.5000 89% 75.8-117 1,4-Dioxane Rejected 0.0250 " 0.5000 81% 54.7-150 cis-1,3-Dichloropropene 0.436 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 " 0.5000 96% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 " <th< th=""><th>trans-1,2-Dichloroethene</th><th>0.501</th><th>J</th><th></th><th>0.0250</th><th>"</th><th>0.5000</th><th></th><th>100%</th><th>0-200</th><th></th><th></th></th<>	trans-1,2-Dichloroethene	0.501	J		0.0250	"	0.5000		100%	0-200			
Cyclohexane 0.441 0.0250 " 0.5000 88% 72.5-119 Benzene 0.494 0.0250 " 0.5000 99% 75.7-118 n-Heptane 0.447 J 0.0250 " 0.5000 89% 66.9-152 1,2-Dichloropropane 0.532 0.0250 " 0.5000 89% 76.8-118 Trichloroethene 0.445 0.0250 " 0.5000 89% 70.1-119 Bromodichloromethane 0.473 0.0250 " 0.5000 89% 75.8-117 1,4-Dioxane Rejected 0.0250 " 0.5000 81% 54.7-150 cis-1,3-Dichloropropene 0.436 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 " 0.5000 72% 62.9-133 trans-1,3-Dichloropropene 0.479 0.0250 " 0.5000 93% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 "	Chloroform	0.491			0.0250	"	0.5000		98%	79.6-115			
Benzene 0.494 0.0250 " 0.5000 99% 75.7-118 n-Heptane 0.447 J 0.0250 " 0.5000 89% 66.9-152 1,2-Dichloropropane 0.532 0.0250 " 0.5000 106% 76.8-118 Trichloroethene 0.445 0.0250 " 0.5000 89% 70.1-119 Bromodichloromethane 0.473 0.0250 " 0.5000 81% 54.7-150 cis-1,3-Dichloropropene 0.436 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 " 0.5000 72% 62.9-133 trans-1,3-Dichloropropene 0.479 0.0250 " 0.5000 96% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 " 0.5000 93% 92.3-106	1,2-Dichloroethane	0.531			0.0250	"	0.5000		106%	79.8-115			
n-Heptane 0.447 J 0.0250 " 0.5000 89% 66.9-152 1,2-Dichloropropane 0.532 0.0250 " 0.5000 106% 76.8-118 Trichloroethene 0.445 0.0250 " 0.5000 89% 70.1-119 Bromodichloromethane 0.473 0.0250 " 0.5000 81% 54.7-150 t-A-Dioxane Rejected 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 " 0.5000 72% 62.9-133 trans-1,3-Dichloropropene 0.479 0.0250 " 0.5000 96% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 " 0.5000 93% 92.3-106	Cyclohexane	0.441			0.0250	"	0.5000		88%	72.5-119			
1,2-Dichloropropane 0.532 0.0250 " 0.5000 106% 76.8-118 Trichloroethene 0.445 0.0250 " 0.5000 89% 70.1-119 Bromodichloromethane 0.473 0.0250 " 0.5000 95% 75.8-117 1,4-Dioxane Rejected 0.0250 " 0.5000 81% 54.7-150 cis-1,3-Dichloropropene 0.436 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 " 0.5000 72% 62.9-133 trans-1,3-Dichloropropene 0.479 0.0250 " 0.5000 96% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 " 0.5000 93% 92.3-106	Benzene	0.494			0.0250	"	0.5000		99%	75.7-118			
Trichloroethene 0.445 0.0250 " 0.5000 89% 70.1-119 Bromodichloromethane 0.473 0.0250 " 0.5000 95% 75.8-117 1,4-Dioxane Rejected 0.0250 " 0.5000 81% 54.7-150 cis-1,3-Dichloropropene 0.436 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 " 0.5000 72% 62.9-133 trans-1,3-Dichloropropene 0.479 0.0250 " 0.5000 96% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 " 0.5000 93% 92.3-106	n-Heptane	0.447	J		0.0250	"	0.5000		89%	66.9-152			
Bromodichloromethane 0.473 0.0250 0.5000 95% 75.8-117 1,4-Dioxane Rejected 0.0250 0.5000 81% 54.7-150 cis-1,3-Dichloropropene 0.436 0.0250 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 0.5000 72% 62.9-133 trans-1,3-Dichloropropene 0.479 0.0250 0.5000 96% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 0.5000 93% 92.3-106	1,2-Dichloropropane	0.532			0.0250	"	0.5000		106%	76.8-118			
1,4-Dioxane Rejected 0.0250 " 0.5000 81% 54.7-150 cis-1,3-Dichloropropene 0.436 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 " 0.5000 72% 62.9-133 trans-1,3-Dichloropropene 0.479 0.0250 " 0.5000 96% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 " 0.5000 93% 92.3-106	Trichloroethene	0.445			0.0250	"	0.5000		89%	70.1-119			
cis-1,3-Dichloropropene 0.436 0.0250 " 0.5000 87% 75.5-115 4-Methyl-2-pentanone Rejected 0.0250 " 0.5000 72% 62.9-133 trans-1,3-Dichloropropene 0.479 0.0250 " 0.5000 96% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 " 0.5000 93% 92.3-106	Bromodichloromethane	0.473			0.0250	"	0.5000		95%	75.8-117			
4-Methyl-2-pentanone Rejected 0.0250 " 0.5000 72% 62.9-133 trans-1,3-Dichloropropene 0.479 0.0250 " 0.5000 96% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 " 0.5000 93% 92.3-106	1,4-Dioxane	Rejected			0.0250	"	0.5000		81%	54.7-150			
trans-1,3-Dichloropropene 0.479 0.0250 " 0.5000 96% 75.8-117 1,1,2-Trichloroethane 0.463 0.0250 " 0.5000 93% 92.3-106	cis-1,3-Dichloropropene	0.436			0.0250	"	0.5000		87%	75.5-115			
1,1,2-Trichloroethane 0.463 0.0250 " 0.5000 93% 92.3-106	4-Methyl-2-pentanone	Rejected			0.0250	"	0.5000		72%	62.9-133			
	trans-1,3-Dichloropropene	0.479			0.0250	"	0.5000		96%	75.8-117			
Dibromochloromethane 0.450 J 0.0250 " 0.5000 90% 69-132	1,1,2-Trichloroethane	0.463			0.0250	"	0.5000		93%	92.3-106			
	Dibromochloromethane	0.450	J		0.0250	"	0.5000		90%	69-132			



LCS (B409106-BS1)

Environmental Protection Agency Region 5

Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605 Phone:(312)353-8370 Fax:(312)886-2591

Air Division, US EPA Region 5 Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Prenared & Analyzed: Sen-19-14

Batch B409106 - ColdTrap Dehydration

LCS (B409106-BS1)				Prepare	a & Anaiy	zea: Sep-1	9-14				
		Flags /		Reporting		Spike	Source		%REC		RPD
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit
Toluene	0.448			0.0250	ppbv	0.5000		90%	73.2-120		
2-Hexanone	Rejected			0.0250	"	0.5000		39%	76-110		
1,2-Dibromoethane (EDB)	0.480			0.0250	"	0.5000		96%	75.5-118		
Tetrachloroethene	0.466			0.0250	"	0.5000		93%	67.1-125		
Chlorobenzene	0.510			0.0250	"	0.5000		102%	68.5-121		
Ethylbenzene	0.468			0.0250	"	0.5000		94%	74.9-118		
m+p-Xylene	0.916			0.100	"	1.000		92%	79.8-121		
Bromoform	0.435			0.100	"	0.5000		87%	72.4-119		
Styrene	0.428			0.100	"	0.5000		86%	71.5-122		
1,1,2,2-Tetrachloroethane	0.490			0.0250	"	0.5000		98%	92-106		
o-Xylene	0.444			0.0500	"	0.5000		89%	77.6-124		
4-ethyltoluene	0.511	J		0.0250	"	0.5000		102%	96.7-122		
1,3,5-Trimethylbenzene	0.513			0.0500	"	0.5000		103%	74.4-121		
1,2,4-Trimethylbenzene	0.492			0.0250	"	0.5000		98%	71.9-126		
1,3-Dichlorobenzene	0.525			0.100	"	0.5000		105%	67.9-132		
Benzyl chloride	0.503			0.0250	"	0.5000		101%	60.7-134		
1,4-Dichlorobenzene	0.521			0.100	"	0.5000		104%	65.4-136		
1,2-Dichlorobenzene	0.517			0.0500	"	0.5000		103%	69.3-129		
1,2,4-Trichlorobenzene	0.506			0.100	"	0.5000		101%	39.7-186		
Naphthalene	0.573	J		0.100	"				40-200		
Surrogate: Dichlorodifluoromethane	0.505				"	0.5000		101%	77.8-116		
Surrogate: Dichlorotetrafluoroethane	0.518				"	0.5000		104%	89-108		
Surrogate: Trichlorofluoromethane	0.486				"	0.5000		97%	78.6-114		
Surrogate:	0.686	J			"	0.5000		137%	75.3-119		
1,1,2-trichloro-1,2,2-trifluoroethane (Freon	0.406				_						
Surrogate: 1,1,1-Trichloroethane	0.496				"	0.5000		99%	92.5-105		
Surrogate: Carbon tetrachloride	0.527				"	0.5000		105%	76.3-118		

		Flags /		Reporting		Spike	Source		%REC		RPD
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit
Propene	0.512			0.0250	ppbv	0.5000		102%	76-112	0.2	19.6
Chloromethane	0.490			0.0250	"	0.5000		98%	77.6-116	3	26.9
Vinyl chloride	0.498			0.0250	"	0.5000		100%	77.4-117	1	25.1



Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605 Phone:(312)353-8370 Fax:(312)886-2591

Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B409106 - ColdTrap Dehydration

LCS Dup (B409106-BSD1)	Prepared & Analyzed: Sep-19-14											
	Flags /		Reporting		Spike	Source		%REC		RPD		
Analyte Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit		
1,3-butadiene 0.412			0.0250	ppbv	0.5000		82%	77.7-115	12	33.2		
Bromomethane 0.536			0.0250	"	0.5000		107%	75.2-117	6	26.6		
Chloroethane 0.584			0.0250	"	0.5000		117%	79.8-114	23	29.5		
Ethanol 0.696	J		0.0500	"	0.5000		139%	50-124	3	200		
Acrolein 0.457			0.0250	"	0.5000		91%	67.9-130	7	29.8		
Isopropyl alcohol 0.334	J		0.0500	"	0.5000		67%	81-106	1	25		
Acetone 0.419	J		0.0250	"	0.5000		84%	60-137	5	28.7		
1,1-Dichloroethene 0.488	J		0.0250	"	0.5000		98%	77.3-118	9	15.9		
Methylene chloride 0.543			0.100	"	0.5000		109%	78.7-117	4	20.7		
Carbon disulfide Rejected			0.0250	"	0.5000		101%	53-140	11	200		
Methyl tert-butyl ether 0.357			0.0250	"	0.5000		71%	79.1-118	14	31.9		
1,1-Dichloroethane 0.404			0.0250	"	0.5000		81%	79.8-117	7	13.1		
Vinyl acetate 0.349			0.0250	"	0.5000		70%	85-106	7	200		
Hexane 0.400			0.0250	"	0.5000		80%	70.1-121	9	43.5		
2-Butanone 0.347			0.0250	"	0.5000		69%	82-108	5	25		
cis-1,2-Dichloroethene 0.418			0.0250	"	0.5000		84%	78.1-115	10	29.6		
trans-1,2-Dichloroethene 0.478	J		0.0250	"	0.5000		96%	0-200	5	25		
Chloroform 0.467			0.0250	"	0.5000		93%	79.6-115	5	25.2		
1,2-Dichloroethane 0.504			0.0250	"	0.5000		101%	79.8-115	5	24.6		
Cyclohexane 0.425			0.0250	"	0.5000		85%	72.5-119	4	34.5		
Benzene 0.451			0.0250	"	0.5000		90%	75.7-118	9	27.4		
n-Heptane 0.417	J		0.0250	"	0.5000		83%	66.9-152	7	25		
1,2-Dichloropropane 0.501			0.0250	"	0.5000		100%	76.8-118	6	25.3		
Trichloroethene 0.427			0.0250	"	0.5000		85%	70.1-119	4	34.1		
Bromodichloromethane 0.443			0.0250	"	0.5000		89%	75.8-117	7	26.5		
1,4-Dioxane Rejected			0.0250	"	0.5000		75%	54.7-150	8	58.6		
cis-1,3-Dichloropropene 0.389			0.0250	"	0.5000		78%	75.5-115	11	31.1		
4-Methyl-2-pentanone Rejected			0.0250	"	0.5000		73%	62.9-133	0.4	42		
trans-1,3-Dichloropropene 0.427			0.0250	"	0.5000		85%	75.8-117	11	31.7		
1,1,2-Trichloroethane 0.446			0.0250	"	0.5000		89%	92.3-106	4	11.5		
Dibromochloromethane 0.432	J		0.0250	"	0.5000		86%	69-132	4	25		
Toluene 0.424			0.0250	"	0.5000		85%	73.2-120	5	30.6		
2-Hexanone Rejected			0.0250	"	0.5000		36%	76-110	7	46.8		
1,2-Dibromoethane (EDB) 0.440			0.0250	"	0.5000		88%	75.5-118	9	31.5		

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Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605 Phone:(312)353-8370 Fax:(312)886-2591

Air Division, US EPA Region 5 Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B409106 - ColdTrap Dehydration

LCS Dup (B409106-BSD1)		Prepared & Analyzed: Sep-19-14										
		Flags /		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	
Tetrachloroethene	0.445			0.0250	ppbv	0.5000		89%	67.1-125	5	13.8	
Chlorobenzene	0.503			0.0250	"	0.5000		101%	68.5-121	1	31.9	
Ethylbenzene	0.464			0.0250	"	0.5000		93%	74.9-118	0.7	31.6	
m+p-Xylene	0.915			0.100	"	1.000		91%	79.8-121	0.1	28.9	
Bromoform	0.426			0.100	"	0.5000		85%	72.4-119	2	34.6	
Styrene	0.429			0.100	"	0.5000		86%	71.5-122	0.2	19.7	
1,1,2,2-Tetrachloroethane	0.476			0.0250	"	0.5000		95%	92-106	3	11.5	
o-Xylene	0.416			0.0500	"	0.5000		83%	77.6-124	7	28.7	
4-ethyltoluene	0.534	J		0.0250	"	0.5000		107%	96.7-122	4	25	
1,3,5-Trimethylbenzene	0.544			0.0500	"	0.5000		109%	74.4-121	6	29.8	
1,2,4-Trimethylbenzene	0.534			0.0250	"	0.5000		107%	71.9-126	8	32.1	
1,3-Dichlorobenzene	0.525			0.100	"	0.5000		105%	67.9-132	0.07	37.9	
Benzyl chloride	0.494			0.0250	"	0.5000		99%	60.7-134	2	48.3	
1,4-Dichlorobenzene	0.520			0.100	"	0.5000		104%	65.4-136	0.1	39.6	
1,2-Dichlorobenzene	0.538			0.0500	"	0.5000		108%	69.3-129	4	34	
1,2,4-Trichlorobenzene	0.529			0.100	"	0.5000		106%	39.7-186	4	77.1	
Naphthalene	0.728	J		0.100	"				40-200	24	200	
Surrogate: Dichlorodifluoromethane	0.532				"	0.5000		106%	77.8-116			
Surrogate: Dichlorotetrafluoroethane	0.536				"	0.5000		107%	89-108			
Surrogate: Trichlorofluoromethane	0.507				"	0.5000		101%	78.6-114			
Surrogate:	0.609	J			"	0.5000		122%	75.3-119			
1,1,2-trichloro-1,2,2-trifluoroethane (Freon	0.477				_							
Surrogate: 1,1,1-Trichloroethane	0.477				"	0.5000		95%	92.5-105			
Surrogate: Carbon tetrachloride	0.498				"	0.5000		100%	76.3-118			

Ouplicate (B409106-DUP1)	Source: 1409007-03	Prepared & Analyzed: Sep-19-14
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		Flags /		Reporting		Spike	Source		%REC		RPD
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit
Propene	0.912	K		0.0250	ppbv		0.903			1	30
Chloromethane	0.462			0.0250	"		0.461			0.2	30
Vinyl chloride	U			0.0250	"		U				30
1,3-butadiene	0.0257	K		0.0250	"		0.0260			1	30
Bromomethane	6.65E-3			0.0250	"		5.68E-3			16	30
Chloroethane	U			0.0250	"		7.73E-3				30

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Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605 Phone:(312)353-8370 Fax:(312)886-2591

Air Division, US EPA Region 5 Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B409106 - ColdTrap Dehydration

Duplicate (B409106-DUP1)	Source: 1409007-03 Prepared & Analyzed: Sep-19-14										
		Flags /		Reporting		Spike	Source		%REC		RPD
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit
Ethanol	2.27	J		0.0500	ppbv		2.25			0.9	40
Acrolein	0.464			0.0250	"		0.437			6	40
Isopropyl alcohol	1.06	J		0.0500	"		1.09			2	40
Acetone	2.63	J		0.0250	"		2.61			0.6	40
1,1-Dichloroethene	U	J		0.0250	"		U				30
Methylene chloride	0.114			0.100	"		0.159			33	30
Carbon disulfide	Rejected			0.0250	"		Rejected			3	40
Methyl tert-butyl ether	U			0.0250	"		U				40
1,1-Dichloroethane	U			0.0250	"		U				30
Vinyl acetate	8.62E-3			0.0250	"		U				40
Hexane	U			0.0250	"		0.366				30
2-Butanone	0.458			0.0250	"		0.462			0.9	50
cis-1,2-Dichloroethene	U			0.0250	"		U				30
trans-1,2-Dichloroethene	5.52E-3	J		0.0250	"		5.89E-3			6	30
Chloroform	0.0272			0.0250	"		0.0266			2	30
1,2-Dichloroethane	9.26E-3			0.0250	"		9.58E-3			3	30
Cyclohexane	0.112			0.0250	"		0.109			3	30
Benzene	0.202			0.0250	"		0.200			1	30
n-Heptane	0.134	J		0.0250	"		0.131			2	30
1,2-Dichloropropane	U			0.0250	"		U				30
Trichloroethene	0.0115			0.0250	"		0.0108			6	30
Bromodichloromethane	U			0.0250	"		U				30
1,4-Dioxane	Rejected			0.0250	"		Rejected				40
cis-1,3-Dichloropropene	U			0.0250	"		U				30
4-Methyl-2-pentanone	Rejected			0.0250	"		Rejected			0.5	40
trans-1,3-Dichloropropene	U			0.0250	"		U				30
1,1,2-Trichloroethane	U			0.0250	"		U				30
Dibromochloromethane	U	J		0.0250	"		U				30
Toluene	0.424			0.0250	"		0.433			2	30
2-Hexanone	Rejected			0.0250	"		Rejected			2	40
1,2-Dibromoethane (EDB)	U			0.0250	"		U				30
Tetrachloroethene	0.0160			0.0250	"		0.0165			3	30
Chlorobenzene	0.0118			0.0250	"		0.0120			2	30
Ethylbenzene	0.0609			0.0250	"		0.0620			2	30



Chicago Regional Laboratory

536 South Clark Street, Chicago, IL 60605 Phone:(312)353-8370 Fax:(312)886-2591

Air Division, US EPA Region 5 Project: BP Whiting Refinery

77 West Jackson BoulevardProject Number: [none]Reported:Chicago IL, 60605Project Manager: Motria CaudillJan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B409106 - ColdTrap Dehydration

Duplicate (B409106-DUP1)	Source:	1409007-03	Prepared & Analyzed: Sep-19-14								
		Flags /		Reporting		Spike	Source		%REC		RPD
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit
m+p-Xylene	0.173			0.100	ppbv		0.177			3	30
Bromoform	8.15E-3			0.100	"		8.25E-3			1	30
Styrene	0.0200			0.100	"		0.0207			3	30
1,1,2,2-Tetrachloroethane	8.56E-3			0.0250	"		U				30
o-Xylene	0.0625			0.0500	"		0.0646			3	30
4-ethyltoluene	0.0791	J		0.0250	"		0.0804			2	30
1,3,5-Trimethylbenzene	0.0252			0.0500	"		0.0254			1	30
1,2,4-Trimethylbenzene	0.112			0.0250	"		0.109			3	30
1,3-Dichlorobenzene	U			0.100	"		U				30
Benzyl chloride	7.41E-3			0.0250	"		0.0252			109	30
1,4-Dichlorobenzene	0.0147			0.100	"		0.0144			2	30
1,2-Dichlorobenzene	U			0.0500	"		U				30
1,2,4-Trichlorobenzene	0.217			0.100	"		0.209			4	30
Naphthalene	1.31	J, K		0.100	"		1.27			4	30
Surrogate: Dichlorodifluoromethane	0.418				"	0.5250		80%	90-120		
Surrogate: Dichlorotetrafluoroethane	0.0188				"	2.000E-2		94%	80-120		
Surrogate: Trichlorofluoromethane	0.204				"	0.2390		85%	90-120		
Surrogate: 1,1,2-trichloro-1,2,2-trifluoroethane (Freon	0.0528	J			"	7.300E-2		72%	90-120		
Surrogate: 1,1,1-Trichloroethane	4.96E-3	J			"	5.000E-3		99%	80-120		
Surrogate: Carbon tetrachloride	0.0692				"	8.600E-2		80%	80-120		

Batch B411044 - ColdTrap Dehydration

Blank (B411044-BLK1)	Prepared & Analyzed: Sep-22-14											
		Flags /		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	
Ethanol	U	J		0.0500	ppbv							
Isopropyl alcohol	0.125	J		0.0500	"							
Acetone	0.106	J		0.0250	"							
Methylene chloride	U			0.100	"							
Surrogate: Dichlorodifluoromethane	U				"	0.5250		%	90-120			
Surrogate: Trichlorofluoromethane	U				"	0.2390		%	90-120			
·												



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Air Division, US EPA Region 5 Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Blank (B411044-BLK1) Analyte	Prepared & Analyzed: Sep-22-14										
		Flags /		Reporting		Spike	Source		%REC		RPD
	Result Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	
Surrogate: Carbon tetrachloride	U				ppbv	8.600E-2		%	80-120		

Blank (B411044-BLK2) Prepared: Sep-22-14 Analyzed: Sep-23-14

Diank (D411044-DERE)	1 repared, Sep-22-14 Amaryzed, Sep-23-14											
		Flags /		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	
Ethanol	U	J		0.0500	ppbv							
Isopropyl alcohol	0.0737	J		0.0500	"							
Acetone	0.110	J		0.0250	"							
Methylene chloride	U			0.100	"							
Surrogate: Dichlorodifluoromethane	U				"	0.5250		%	90-120			
Surrogate: Trichlorofluoromethane	U				"	0.2390		%	90-120			
Surrogate: Carbon tetrachloride	U				"	8.600E-2		%	80-120			

LCS (B411044-BS1) Prepared & Analyzed: Sep-22-14

ECS (BTITOTT BST)	1 repared & Amaryzed, Sep-22-14										
		Flags /		Reporting		Spike	Source		%REC		RPD
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit
Ethanol	0.628	J		0.0500	ppbv	0.5000		126%	50-124		
Isopropyl alcohol	0.400	J		0.0500	"	0.5000		80%	81-106		
Acetone	0.436	J		0.0250	"	0.5000		87%	60-137		
Methylene chloride	0.513			0.100	"	0.5000		103%	78.7-117		
Surrogate: Dichlorodifluoromethane	0.475				"	0.5000		95%	77.8-116		
Surrogate: Trichlorofluoromethane	0.504				"	0.5000		101%	78.6-114		
Surrogate: Carbon tetrachloride	0.511				"	0.5000		102%	76.3-118		

LCS Dup (B411044-BSD1) Prepared & Analyzed: Sep-22-14

Ees Dup (Billivii BSD1)	Trepured & Thursteen Sep 22 11											
		Flags /		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	
Ethanol	0.674	J		0.0500	ppbv	0.5000		135%	50-124	7	200	
Isopropyl alcohol	0.355	J		0.0500	"	0.5000		71%	81-106	12	25	
Acetone	0.430	J		0.0250	"	0.5000		86%	60-137	1	28.7	
Methylene chloride	0.524			0.100	"	0.5000		105%	78.7-117	2	20.7	
Surrogate: Dichlorodifluoromethane	0.486				"	0.5000		97%	77.8-116			

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Ethanol

Acetone

Isopropyl alcohol

Methylene chloride

Surrogate: Dichlorodifluoromethane

Surrogate: Trichlorofluoromethane

Surrogate: Carbon tetrachloride

Environmental Protection Agency Region 5

Chicago Regional Laboratory

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Air Division, US EPA Region 5 Project: BP Whiting Refinery

77 West Jackson Boulevard Project Number: [none] Reported:
Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B411044 - ColdTrap Dehy	dration											
LCS Dup (B411044-BSD1)	Prepared: Sep-22-14 Analyzed: Sep-23-14											
		Flags /		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	
Surrogate: Trichlorofluoromethane	0.501				ppbv	0.5000		100%	78.6-114			
Surrogate: Carbon tetrachloride	0.515				"	0.5000		103%	76.3-118			
LCS Dup (B411044-BSD2)				Prepared: Se	ep-22-14 <i>A</i>	analyzed: S	ер-23-14					
		Flags /		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	
Ethanol	0.612	J		0.0500	ppbv	0.5000		122%	50-124	200	200	
Isopropyl alcohol	0.355	J		0.0500	"	0.5000		71%	81-106	200	25	
Acetone	0.435	J		0.0250	"	0.5000		87%	60-137	200	28.7	
Methylene chloride	0.523			0.100	"	0.5000		105%	78.7-117	200	20.7	
Surrogate: Dichlorodifluoromethane	0.490				"	0.5000		98%	77.8-116			
Surrogate: Trichlorofluoromethane	0.446				"	0.5000		89%	78.6-114			
Surrogate: Carbon tetrachloride	0.532				"	0.5000		106%	76.3-118			
Duplicate (B411044-DUP1)	Source: 1	409007-03RE1	l	Prepare								
		Flags /		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	
Ethanol	2.42	J		0.500	ppbv		2.26			7	40	
Acetone	1.80	J		0.250	"		2.80			43	40	
Surrogate: Dichlorodifluoromethane	0.450				"	0.5250		86%	90-120			
Surrogate: Trichlorofluoromethane	0.190				"	0.2390		79%	90-120			
Duplicate (B411044-DUP2)	Source: 1	409007-04RE1		Prepared: S	ep-22-14 <i>A</i>	analyzed: S	ер-23-14					
		Flags /		Reporting		Spike	Source		%REC		RPD	
Analyte	Result	Qualifiers	MDL	Limit	Units	Level	Result	%REC	Limits	RPD	Limit	

0.0500

0.0500

0.0250

0.100

ppbv

2.33

1.02

2.42

0.164

0.5250

0.2390

8.600E-2

J

J

2.29

1.01

2.46

0.175

0.413

0.196

0.0664

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40

40

40

30

Report Name: 1409007 FINAL Jan 13 15 1458

90-120

90-120

80-120

79%

82%

77%

2

0.5

2



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77 West Jackson Boulevard Project Number: [none]

Chicago IL, 60605 Project Manager: Motria Caudill

Reported:Jan-13-15 14:58

Air Toxics by GC/MS, EPA TO-15 (modified) - Quality Control US EPA Region 5 Chicago Regional Laboratory

Batch B411044 - ColdTrap Dehydration

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Chicago Regional Laboratory

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Air Division, US EPA Region 5 Project: BP Whiting Refinery

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Chicago IL, 60605 Project Manager: Motria Caudill Jan-13-15 14:58

Notes and Definitions

R Rejected

K The identification of the analyte is acceptable; the reported value may be biased high. The actual value is expected to be less than

the reported value.

J The identification of the analyte is acceptable; the reported value is an estimate.

U Not Detected

NR Not Reported

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